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## 1. INTRODUCTION

The National Centers for Environmental Prediction (NCEP) regional spectral model (RSM) developed by Juang and Kanamitsu (1994) is a limited area spectral model for regional weather and climate prediction and operationally being used as a member of short-range ensemble forecast system at NCEP. The RSM was designed to have the same model structure, dynamics, and physics as the NCEP global forecast system (GFS) that is a global spectral model for global weather and climate prediction. While in the GFS, total fields are predicted using spherical harmonic functions for spectral computation, in the RSM, perturbation fields from coarse-resolution base fields (usually global analysis or forecast fields) are represented by double Fourier trigonometric functions, which are relaxed to zero at lateral boundaries. As a significant improvement in the GFS physics has been accomplished, we have implemented the updated GFS physics into the RSM recently. More details of the updated physics are described in the following section.

The efforts for improving weather forecast have drawn the model to be run at higher resolution with more sophisticated physics, which requires higher model performance in computation speed. Furthermore, recent development of ensemble forecast system and extensive activities in regional climate simulations using limited area model needed a significant improvement in model performance. Toward this effort, most of up-to-date operational weather forecast models take advantage of massive parallel processor computer to improve their performance. The parallel version of the old RSM at NCEP (hereafter old RSM-MPI) is based on 1D decomposition using message passing interface (MPI) which enables data exchange among distributed memories. However, the performance of the old RSM-MPI has been found to fall short of NCEP's current demands for the model to be run at higher resolution (e.g., about 30km over North America). To improve the RSM computing performance further, we have developed a fully parallelized RSM-MPI code based on 2D

decomposition on IBM-SP, a distributed-memory parallel supercomputer. Efforts for additional performance improvement are also presented in this paper.

## 2. MODEL PHYSICS UPDATE

The RSM model physics has been successfully updated consistent with recent update of the GFS model physics, which was written on pressure coordinate system rather than on the sigma coordinate system. In the update, the diagnostic large-scale cloud scheme is replaced by a prognostic cloud scheme. Radiation parameterization uses the predicted cloud condensate in the cloud-radiation interaction. The Rapid Radiative Transfer Model (RRTM) is now used for the longwave radiation. A new prognostic variable Ozon is introduced, which is predicted using temporally and spatially varying climatological Ozon production and destruction rate. For the cumulus convection that uses a simplified Arakawa-Schubert scheme, mass fluxes induced in the updraft and the downdraft are now allowed to transport momentum, which has produced a significant reduction of the false alarm rate for tropical storm forecasts. The gravity wave drag parameterization includes the effects of subgrid orographic asymmetry and fractional area of subgrid orography larger than grid orographic height for 4 different wind directions in addition to subgrid orographic variance in old version. The Navy 10 minute topography data is replaced by high resolution USGS topography data in which 2 minute, 4 minute, and 8 minute resolution data are available. The surface soil and vegetation types are changed to have heterogeneous types with 9 and 13 kinds, respectively. Surface albedo now has 4 components. More details of the updated model physics can be found in Moorthi et al. (2001) and Hou et al. (2002).

## 3. FULL PARALLELIZATION OF RSM WITH MPI

A spectral model essentially requires a transformation from spectral to grid-point space or vice versa using the Fast Fourier Transform (FFT) algorithm. Because the FFT algorithm requires all of grid-point values within the row or column, 3D decomposition for parallelization in a spectral model

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is almost impossible while 1D or 2D decomposition is possible.

The old RSM-MPI is based on 1D decomposition using MPI. The disadvantage of 1D decomposition is that the number of working processors is limited by the vertical layers which are usually much smaller than horizontal grid points in NWP models. In the old RSM-MPI, furthermore, the number of vertical layers assigned in each processor is designed to be the same and thus, it has to be a factor of the number of the whole vertical layers.

The new 2D decomposition which we developed in this study has the advantage to use more processors than the old 1D decomposition because there are much more grid points in 2D to be partitioned. Figure 1 shows the schematic diagram for 2D decomposition used in the new RSM-MPI code with an example of 12 processors. Each stage is characterized by full dimension in one of three directions. As mentioned earlier, in a spectral model a full dimension in one of two horizontal directions (East-West [X] and North-South [Y] directions in Fig. 1) is needed for the FFT computation. Similar method has been introduced in several previous studies (Oikawa, 2001; Juang et al., 2001; Foster and Worley, 1997; Barros et al., 1995; Skalin and Borge, 1997). First, rows along X and Y directions in the spectral space are divided depending on the number of working processors while the vertical column has a full dimension (upper left corner cube in Fig. 1). Then, for transformation from spectral to grid-point space in Y direction using FFT algorithm, rows along X direction and vertical columns are divided while Y direction has a full dimension (upper middle and right corner cubes in Fig. 1), for which the transpose (rearrangement of partial and full dimensions as shown in upper left corner and middle cubes in Fig. 1) is performed using the MPI routines. Then, for transformation from spectral to grid-point space in X direction, rows along Y direction and vertical columns are divided while X direction has a full dimension (lower right corner and middle cubes in Fig. 1), again using the MPI routines for the transpose (upper and lower right corner cubes in Fig. 1). After the last transpose (lower middle and left corner cubes in Fig. 1), we have partial dimensions along the horizontal directions but a full dimension along the vertical direction in grid-point space (lower left corner cube in Fig. 1) where nonlinear model dynamics and model physics that entails vertical data dependency are computed. The transformation back to the spectral space is conducted in the reverse way denoted as reverse arrows in the Fig. 1. The computation of the semi-implicit time integration is conducted in the spectral space with partial dimensions along the horizontal directions but a full dimension along the vertical direction (upper left corner cube in Fig. 1).

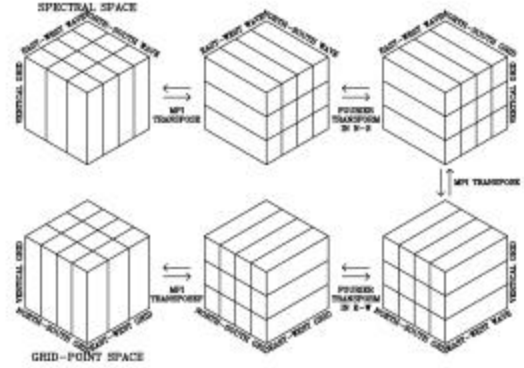


Figure 1. Schematic diagram of 2D decomposition with 12 processors for transformation between spectral to grid-points space using MPI transpose.

The performance of the model with MPI can be represented in terms of speedup. The speedup,  $S_p(n)$ , is defined as the ratio of the time spent by single processor run and the one by  $n$  processor run. For the perfect parallelization, therefore,  $S_p(n)=n$ . Practically, however,  $S_p(n)$  is always less than  $n$  because there are not only sequential codes that cannot be parallelized but also there is a time spent for communication. Assuming that the time spent in communication is negligible, the speedup can be expressed as (e.g., see Juang et al., 2003)

$$S_p(n) = \frac{s + p}{s + p/n} \quad (1)$$

where  $s$  is the time spent by the sequential portion of the code in the computation and  $p$  the time spent by the parallel portion of the code in computation. Therefore, Eq. (1) indicates a maximal speedup achievable theoretically for a percentage of parallelization for given  $n$  processors. This theoretical speedup curve is used to evaluate the speedup performance of the new RSM-MPI code.

The speedup performance of the 24-hour RSM runs at NCEP IBM-SP machine with 4, 8, 16, 32, 42, and 64 working processors is shown in Fig. 2 for two different dimensions. Note that the results for the old RSM-MPI with processors more than 42 are not available because the number of working processors in the old RSM-MPI code is limited by the number of the vertical grid points. For the smaller dimension of  $(X,Y,Z)=(120,121,42)$  (Fig. 2a), the speedup of the new RSM-MPI is in between those of the theoretically maximal values [i.e., Eq. (1)] achievable from 95% and 99% of parallelization, while the old RSM-MPI shows less speedup than the theoretical one from 95% of parallelization. For the larger dimension of  $(X,Y,Z)=(240,241,42)$  (Fig. 2b), the speedup of the new RSM-MPI is about same as the theoretical one achievable from 99% of parallelization. Compared to the old RSM-MPI, overall speedup performance of the new RSM-MPI is about twice larger.

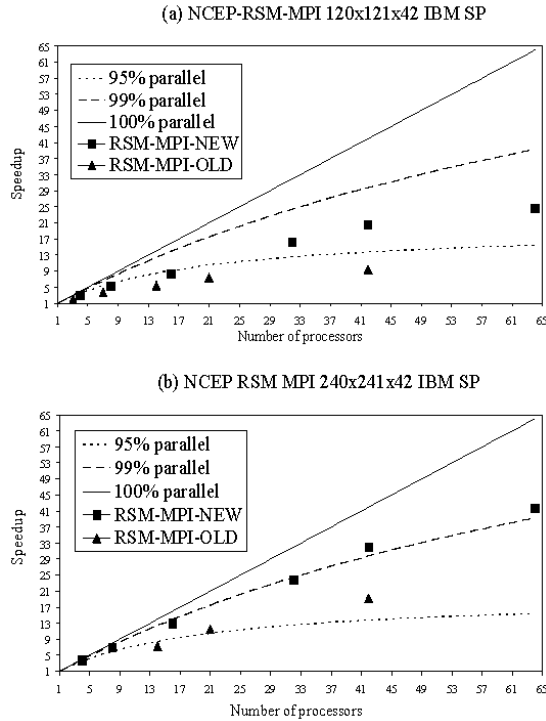


Figure 2. Performance of parallel RSM.

In the MPI code, the memory usage generally decreases with increasing working processors because each processor has smaller dimension with increasing working processors. Note that for the larger dimension (Fig. 2b), the runs of the old RSM-MPI with 4 and 8 working processors were failed because of too much large dimension associated with smaller number of working processors. However, the runs with the new RSM-MPI were successful even for much larger dimension. This is because for the same working processors, each processor has smaller dimension in 2D decomposition than in 1D decomposition. Although not shown in a figure, our tests showed that for a limited computer resource, the new RSM-MPI could use about factor of 5 larger dimensions than the old RSM-MPI, indicating that the memory usage is much smaller in the new RSM-MPI than in the old RSM-MPI.

#### 4. FURTHER PERFORMANCE IMPROVEMENT

The current RSM-MPI uses 8-byte floating point for real variables. A performance test was conducted using 4-byte floating point only for MPI communication, which gave rise to about 20% speed-up without noticeable change in the results. The RSM performance was further improved by correcting indexing in FORTRAN do-loops for some source files, which gave rise to about 20% speed-up. Examples of bad and good indexing on IBM-SP machine are:

##### Bad indexing

```
DO I=1,NX
  DO J=1,NX
    A(I,J)=B(I,J)+C(I,J)
  ENDDO
ENDDO
```

##### Good indexing

```
DO J=1,NX
  DO I=1,NX
    A(I,J)=B(I,J)+C(I,J)
  ENDDO
ENDDO
```

To improve model performance further, parallelization in model input/output is being developed.

#### 5. SUMMARY

The improved GFS model physics has been implemented into the RSM. The updates in radiation, grid-scale condensation, cumulus convection, and gravity wave drag are the most noticeable. A new fully parallelized RSM using MPI based on 2D decomposition has been also developed, which shows much improvement in speedup and memory usage compared to the previous RSM-MPI that is based on 1D decomposition. While in the old RSM-MPI the number of working processors is limited by vertical layers, in the new RSM-MPI it is limited by horizontal dimension, which is usually much larger than vertical dimension in numerical weather models. Compared to the old RSM-MPI, the new RSM-MPI is about twice faster and can use about factor of 5 larger dimensions. The RSM performance was further improved by using 4-byte floating point rather than 8-byte floating point for real variables only in MPI communication and by correcting indexing in FORTRAN do-loops, of which the combined effects gave rise to about 40% speed-up.

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